From: Benjamin.Shorr@noaa.gov Eric Blischke/R10/USEPA/US@EPA To: Subject: Re: A few modifications to the table

Date: 01/11/2007 01:08 PM

Attachments: cumulative distribution 20071008.txt

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Eric-
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Attached is a macro in Excel that creates the fields for the cumulative distribution in a text file- I hope either Margaret or Jim will know how to use it. Either way, I'm working on finishing some code that will generate the graphs as well for the distribution & also mean segments (RM, F&T). Also coded the spatial join in ArcView- need to modify so can do a batch of contaminants at once.

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---- Original Message ----
From: Blischke.Eric@epamail.epa.gov
Date: Thursday, January 11, 2007 3:00 pm
Subject: Re: A few modifications to the table
> Ben, I don't think I received the pdf.
> Eric
                          Benjamin.Shorr@n
                          oaa.gov
```

То

01/11/2007 12:56 Margaret Spence

ΡМ <mspence@parametrix.com>

CC

> Dana

Jay.Field@noaa.gov, Robert

Gensemer

<rgensemer@parametrix.com>,

Davoli/R10/USEPA/US@EPA, Eric

Blischke/R10/USEPA/US@EPA,

Jay.Field@noaa.gov,

Robert.Neely@noaa.gov, Carrie

Smith

> <csmith@parametrix.com>, Jim Koloszar

<jkoloszar@parametrix.com>

> Subject

> the

Re: A few modifications to

table

Thanks guys-

> Attached is a pdf with an example (PAH) of the analyses that I've been > doing for surface sediment. This is for ecorisk so far... I've done > pieces for most other contaminants and have created some macros for > thecumulative distribution, graphing and also for the spatial join.

> B

> ---- Original Message ----> From: Margaret Spence <mspence@parametrix.com> > Date: Thursday, January 11, 2007 1:39 pm

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> Subject: Re: A few modifications to the table
  > Hi gang. I'm working at home again today and will be joining in the > conference call. I've attached an Excel spreadsheet (and a PDF > print-out of it) I worked up yesterday to guide me through the > analysisprocess and keep files, etc. organized. It also includes > severalquestions I came up with yesterday as I started working
      Hopefully, these can get resolved during today's call.
  > If anybody needs to reach me, call my cell phone at (b) (6)
> > Ben, (b) (6)
      Margaret
      Margaret Spence
Phone: 425-458-6369
Fax: 425-458-6363
      mspence@parametrix.com
      PARAMETRIX
      Inspired people - Inspired solutions - Making a difference
      >>> <Benjamin.Shorr@noaa.gov> 01/10/07 10:31 PM >>>
      Sounds good-
      I am trying to use the spreadsheet as a guide- added a couple of
> > fieldsfor Cumulative Distr. charts, summary by areas graphs & maps
  > checking them off as I go...
      (b) (6)
            Perhaps tomorrow after the call and discussion of progress we
> > can see
>> if I need to spread some pieces that I am responsible for to
> Margaret,> Jim or Carrie to ensure that they get enough attention.
      Thanks,
          --- Original Message ----
>> From: Robert Gensemer <rgensemer@parametrix.com>
>> Date: Wednesday, January 10, 2007 6:14 pm
>> Subject: Re: A few modifications to the table
         I think we need to be as consistent with QM as possible in
> terms of
> > > numbers and units. Lets not get too concerned about cleaning up > > every> aspect of the risk parameters table to be a perfect match > > with QM,
   > > though. Remember this is a guide of analyses to do and a
> > compliation of
>>> screening values, not necessarily a formal spreadsheet work
> template> > (unless you guys have decided to do so??). Thanks to all,
      > -Bob
  > > Robert W. Gensemer, Ph.D.
> > Parametrix, Inc.
> > 33972 Texas Street SW
     > Albany, OR 97321
> T 541-791-1667, x-6510
> F 541-791-1699
> C 541-760-1511
      > rgensemer@parametrix.com
>>> > Renjamin.Shorr@noaa.gov> 1/10/2007 7:36:38 AM >>>
> > > Eric-
>>> A few notes on the surface sediment screening numbers for > ecological>> risk:
         I strongly recommend that the units that are in this
> spreadsheet be
> > changed to reflect the units in Query Manager. There should be a
> > Changed to Telesco ....
> > column
> > with the units for each analyte (most metals in PPM, vols/svols
> etc> > PPB), and the guidelines should be adjusted to that for
> consistency.>
  consistency.> >
> Total PCB's TEC should probably be .0598 (off by 10^3)
      > Dieldrin (PPB) numbers are TEC/PEC = 1.9/61.8; spreadsheet has
   > > 2.85/6.7
      ^{>} 2378 TCDD- there is one sample over 9 ng/kg (9E^-6 mg/kg) at 111 under> railroad bridge. Looking directly at TCDD2378 conc. may benefit from
      > paired number.
      > Hexachlorocyclohexane differs from QM TEC/PEC which is 2.37/4.99 PPB,> spreadsheet has .94/1.38
  >> Hexachlorobutadiene, Tetrachloroethene, Trichloroethene units
> may be > > in
         incorrect in spreadsheet (off by 10^3)
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> > Please let me know if there is a call today that I can join- > otherwise> I'm available for the 1pm call tomorrow.
> > > Ren
     > ---- Original Message ----
> From: Blischke.Eric@epamail.epa.gov
> Date: Tuesday, January 9, 2007 3:05 pm
> Subject: Re: A few modifications to the table
     > > Dana, here is a response to your questions and modifications to the> > table. I am copying the data evaluation folks and attaching your> > modifications to the table. I also have a few questions for Ben
     > > regarding how QM handles certain summed values.
       > I do not want to look at aluminum. 7600 mg/kg while screening
  > in
     >> HQ of 0.1 is probably below background - upstream aluminum
>> concentrations range from 12,000 - 33,000 mg/kg. Further, the
     > direct
        > contact exposure scenarios are very conservative (350 days a
     year
> > for a
> > beach?).
  >
       > Regarding the TEQs and DDT, DDE and DDD sums - by manually, I
     > meant
     >> thatit was not being calculated automatically by Query Manager.
     > > shouldbe able to do this in excel. I certainly support looking
  >
     > at
     > the TEQs > but I want to get started on some easier evaluations first. We
     > > haveto prioritize things here.
     >> Ben: What is included in the reported TEQ value - dioxin TEQs
  > or
  > > dioxinand dioxin-like PCB TEQ?
  > > > I don't really know how to best evaluate the PAHs. Reg. > > naphthalene and Benzo(a)pyrene, we can look at these as individual> > > chemicals. Hopefully, if we look at total PAHs,
> total low
     molecular> > weight PAHs and BAP and naphthalene, we will get a
>> sense of the PAH
>> > distribution to help us focus our evaluation. Another thing we
> > > might
 > > want to do is query the carcinogenic PAHs and look at total > > carcinogenicPAHs screened against BAP screening numbers.
         Ben: Do you know high molecular weight and low molecular weight
     > PAHs
        > are calculated.
     >> Regarding the modified table. I am ok with screening non-
     >> carcinogens at
>> 0.1 (with the exception of Aluminum). Because QM is good at
       > looking at
       > concentration ranges, we should look at both HQ = 1 and HQ =
  0.1.>>>
>>> I noticed the error regarding the residential soil PRG for BAP
     > (units
     > > problem). You have correctly modified the screening number to
     be
       > 0.062 mg/kg.
       > Lets figure out the best way to too look at total PCBs (total
  >
     > aroclors
     > or total congeners). For surface water, we should look at total > congeners due to interferences associated with the aroclor
     > results.
     > > For
     > > sediment, we should look at both total congeners and total arocIors.> > The total congeners represents a better number. However, we have
  >
     > much
     > less congener data than aroclor data. (PMX and Ben, I am > > attaching a > > write up on summing).
     >> Regarding TBT in Fish, our TBT data is limited to clams, and
     > juvenile
          Chinook. Only one sample (a clam sample from the shipyard)
     > exceeds
          thefish screening value (detected concentration = 530 ug/kg;
  > fish
     > > screeningnumber = 144 ug/kg; shellfish screening number = 1170 > > ug/kg). We can > > still look at TBT in surface water.
       > Eric
     > > (See attached file: 20070108Davoli Modif to ERIC
     > > RiskParameters.xls)(Seeattached file: 20060201 Kissinger
> > Approach
> > Approacn
> > > Portland Harbor Upstream Fish
> > > Tissue Sample Total PCBs, PCB TEQs, Dioxin_Furan TEQs.doc)
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danadavoli
                                                                  <danadavoli@avva
                                                                 nta.com>
             > > To
                                                                                                                                                  Eric
            > > Blischke/R10/USEPA/US@EPA
> > 01/08/2007 09:44
     >
            > > CC
> >
                                                                                                                                                 Dana
      > Davoli/R10/USEPA/US@EPA
                   > Subject
                                                                                                                                                 A few modifications to the
             > > table
            > >
> >
            > >
             >>> > > > I just checked the HH table. Changes are in yellow.
             > The major changes are HQ=1.0 to HQ=0.1 for the direct
> contact. I
> > > added
> > > AL back in for the beaches because it screens in at HQ = 0.1. I
> > > don't
^{\circ} > > > have the LWG website so I couldn't check if AL screens in for > the> > > in-water sediments.
            >\, I think we only have Aroclors for the beaches, not congeners. I >\, startedto add all of the TEQs that I would like to see (d/f,
            >> started to the sum
>> of these) to the lists but decided to wait until we talk. I
>> don't
>> > thinkit would be that hard for Parametrix to do the
             calculations
> > > in
> > > EXECUTE STATES 
>>>> fromcongeners and the DDEs, DDDs, and DDTs.None of this > should be
> > > done
> > > manually.
> > > >
                   > For PAH, I do not know how the NOAA database defines hi MW
    > versus
> > low MW
> > PAHs so I can't tell how close the hi MW would be to the
> carcinogenic
> > PAHs (B(a)P equivalents.)
>>> > Wasn't sure what you meant by using naphthalene and B(a)P as >>> surrogates.For example, do you mean using he naphthalene tox
>> > > Surrogates::
>> > values
>> > as surrogates
```